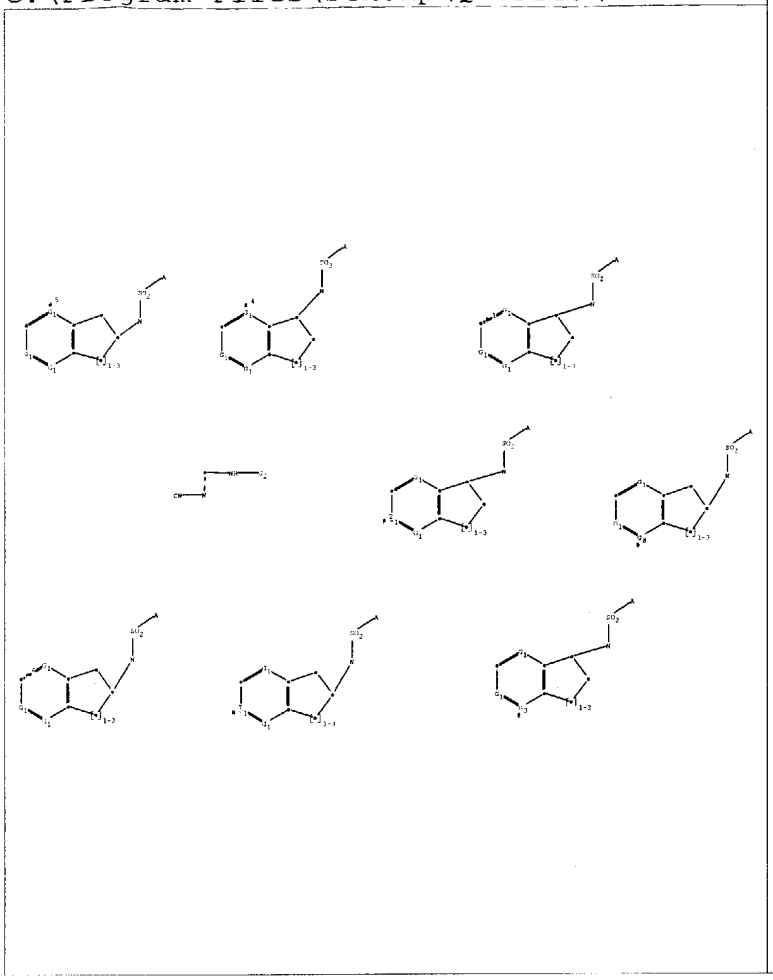
C:\Program Files\Stnexp\Queries\10823987



chain nodes :
14 15 17 27 28 39 40 51 52 60 61 62 76 77 91 92 106 107
121 122 128
ring nodes :
1 2 3 4 5 6 7 8 9 18 19 20 21 22 23 24 25 26 30 31 32
33 34 35 36 37 38 42 43 44 45 46 47 48 49 50 67 68 69 70
71 72 73 74 75 82 83 84 85 86 87 88 89 90 97 98 99 100
101 102 103 104 105 112 113 114 115 116 117 118 119 120
ring/chain nodes :
16 29 41 53 78 93 108 123
chain bonds :
7-14 14-15 15-16 17-60 17-128 24-27 27-28 28-29 36-39 39-40 40-41
48-51 51-52 52-53 60-61 61-62 74-76 76-77 77-78 89-91 91-92 92-93
104-106 106-107 107-108 119-121 121-122 122-123
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 18-19 18-23 19-20
20-21 21-22 22-23 22-24 23-26 24-25 25-26 30-31 30-35 31-32 32-33
33-34 34-35 34-36 35-38 36-37 37-38 42-43 42-47 43-44 44-45 45-46
46-47 46-48 47-50 48-49 49-50 67-68 67-72 68-69 69-70 70-71 71-72
71-73 72-75 73-74 74-75 82-83 82-87 83-84 84-85 85-86 86-87 86-88
87-90 88-89 89-90 97-98 97-102 98-99 99-100 100-101 101-102
101-103 102-105 103-104 104-105 112-113 112-117 113-114 114-115
115-116 116-117 116-118 117-120 118-119 119-120
exact/norm bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 7-14 8-9 14-15 15-16 17-60
17-128 18-19 18-23 19-20 20-21 21-22 22-23 22-24 23-26 24-25
24-27 25-26 27-28 28-29 30-31 30-35 31-32 32-33 33-34 34-35 34-36
35-38 36-37 36-39 37-38 39-40 40-41 42-43 42-47 43-44 44-45 45-46
46-47 46-48 47-50 48-49 48-51 49-50 51-52 52-53 60-61 61-62 67-68
67-72 68-69 69-70 70-71 71-72 71-73 72-75 73-74 74-75 74-76 76-77
77-78 82-83 82-87 83-84 84-85 85-86 86-87 86-88 87-90 88-89 89-90
89-91 91-92 92-93 97-98 97-102 98-99 99-100 100-101 101-102
101-103 102-105 103-104 104-105 104-106 106-107 107-108 112-113
112-117 113-114 114-115 115-116 116-117 116-118 117-120 118-119
119-120

119-121 121-122 122-123

G1:C,N

G2:[*1],[*2],[*3],[*4],[*5],[*6],[*7],[*8]

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom
14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:Atom 19:Atom 20:Atom 21:Atom
22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:CLASS 28:CLASS 29:CLASS
30:Atom 31:Atom 32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom
38:Atom 39:CLASS 40:CLASS 41:CLASS 42:Atom 43:Atom 44:Atom 45:Atom
46:Atom 47:Atom 48:Atom 49:Atom 50:Atom 51:CLASS 52:CLASS 53:CLASS
60:CLASS 61:CLASS 62:CLASS 67:Atom 68:Atom 69:Atom 70:Atom 71:Atom
72:Atom 73:Atom 74:Atom 75:Atom 76:CLASS 77:CLASS 78:CLASS 82:Atom
83:Atom 84:Atom 85:Atom 86:Atom 87:Atom 88:Atom 89:Atom 90:Atom
91:CLASS 92:CLASS 93:CLASS 97:Atom 98:Atom 99:Atom 100:Atom 101:Atom
102:Atom 103:Atom 104:Atom 105:Atom 106:CLASS 107:CLASS 108:CLASS
112:Atom 113:Atom 114:Atom 115:Atom 116:Atom 117:Atom 118:Atom
119:Atom 120:Atom 121:CLASS 122:CLASS 123:CLASS 128:CLASS

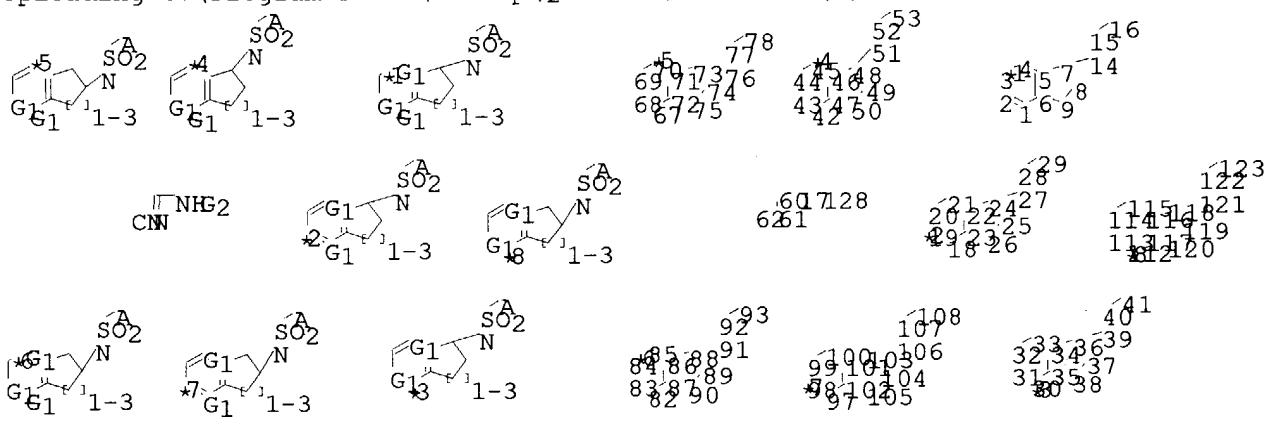
=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=> screen 2016 OR 2026 OR 2039 OR 2040 OR 2045 OR 2047

L1 SCREEN CREATED

=>
Uploading C:\Program Files\Stnexp\Queries\10823987 (a).str



chain nodes :
14 15 17 27 28 39 40 51 52 60 61 62 76 77 91 92 106 107 121 122 128
ring nodes :
1 2 3 4 5 6 7 8 9 18 19 20 21 22 23 24 25 26 30 31 32 33 34 35 36 37 38 42 43 44 45 46 47 48 49 50 67 68 69 70 71 72 73 74 75 82 83 84 85 86 87 88 89 90 97 98 99 100 101 102 103 104 105 112 113 114 115 116 117 118 119 120
ring/chain nodes :
16 29 41 53 78 93 108 123
chain bonds :
7-14 14-15 15-16 17-60 17-128 24-27 27-28 28-29 36-39 39-40 40-41 48-51 51-52 52-53 60-61 61-62 74-76 76-77 77-78 89-91 91-92 92-93 104-106 106-107 107-108 119-121 121-122 122-123

ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 18-19 18-23 19-20 20-21 21-22
22-23 22-24 23-26 24-25 25-26 30-31 30-35 31-32 32-33 33-34 34-35 34-36
35-38 36-37 37-38 42-43 42-47 43-44 44-45 45-46 46-47 46-48 47-50 48-49
49-50 67-68 67-72 68-69 69-70 70-71 71-72 71-73 72-75 73-74 74-75 82-83
82-87 83-84 84-85 85-86 86-87 86-88 87-90 88-89 89-90 97-98 97-102
98-99 99-100 100-101 101-102 101-103 102-105 103-104 104-105 112-113
112-117 113-114 114-115 115-116 116-117 116-118 117-120 118-119 119-120
exact/norm bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 7-14 8-9 14-15 15-16 17-60 17-128
18-19 18-23 19-20 20-21 21-22 22-23 22-24 23-26 24-25 24-27 25-26 27-28
28-29 30-31 30-35 31-32 32-33 33-34 34-35 34-36 35-38 36-37 36-39 37-38
39-40 40-41 42-43 42-47 43-44 44-45 45-46 46-47 46-48 47-50 48-49 48-51
49-50 51-52 52-53 60-61 61-62 67-68 67-72 68-69 69-70 70-71 71-72 71-73
72-75 73-74 74-75 74-76 76-77 77-78 82-83 82-87 83-84 84-85 85-86 86-87
86-88 87-90 88-89 89-90 89-91 91-92 92-93 97-98 97-102 98-99 99-100
100-101 101-102 101-103 102-105 103-104 104-105 104-106 106-107 107-108
112-113 112-117 113-114 114-115 115-116 116-117 116-118 117-120 118-119
119-120 119-121 121-122 122-123

G1:C,N

G2:[*1],[*2],[*3],[*4],[*5],[*6],[*7],[*8]

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 14:CLASS
15:CLASS 16:CLASS 17:CLASS 18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom
24:Atom 25:Atom 26:Atom 27:CLASS 28:CLASS 29:CLASS 30:Atom 31:Atom 32:Atom
33:Atom 34:Atom 35:Atom 36:Atom 37:Atom 38:Atom 39:CLASS 40:CLASS 41:CLASS
42:Atom 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:Atom 49:Atom 50:Atom
51:CLASS 52:CLASS 53:CLASS 60:CLASS 61:CLASS 62:CLASS 67:Atom 68:Atom
69:Atom 70:Atom 71:Atom 72:Atom 73:Atom 74:Atom 75:Atom 76:CLASS 77:CLASS
78:CLASS 82:Atom 83:Atom 84:Atom 85:Atom 86:Atom 87:Atom 88:Atom 89:Atom
90:Atom 91:CLASS 92:CLASS 93:CLASS 97:Atom 98:Atom 99:Atom 100:Atom
101:Atom 102:Atom 103:Atom 104:Atom 105:Atom 106:CLASS 107:CLASS 108:CLASS
112:Atom 113:Atom 114:Atom 115:Atom 116:Atom 117:Atom 118:Atom 119:Atom
120:Atom 121:CLASS 122:CLASS 123:CLASS 128:CLASS

L2 STRUCTURE UPLOADED

=> que L2 NOT L1

L3 QUE L2 NOT L1

=> d l3
L3 HAS NO ANSWERS
L1 SCR 2016 OR 2026 OR 2039 OR 2040 OR 2045 OR 2047
L2 STR
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.
L3 QUE L2 NOT L1

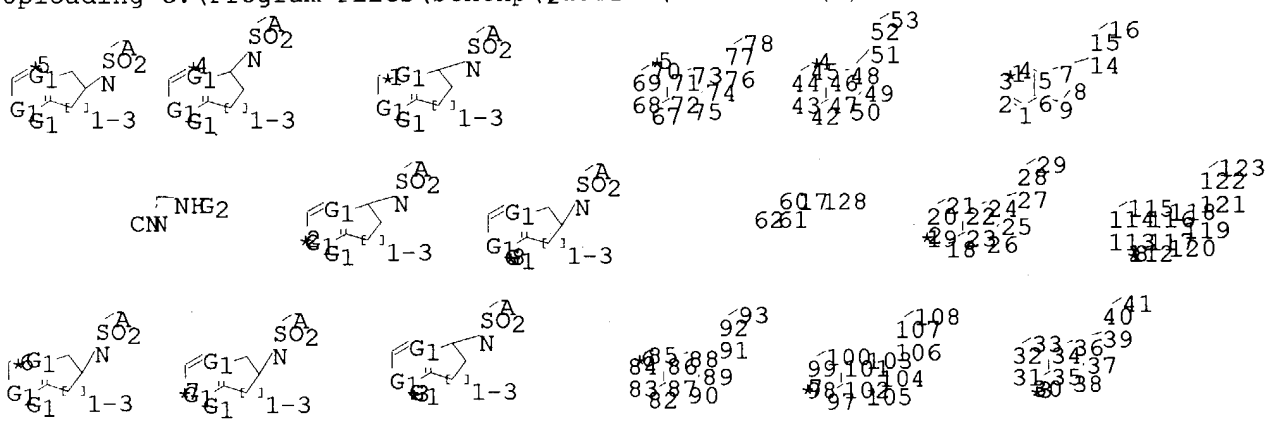
```
=> s l3 sss sam
SAMPLE SEARCH INITIATED 17:49:58 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -      9 TO ITERATE

100.0% PROCESSED       9 ITERATIONS                                0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:             9 TO      360
PROJECTED ANSWERS:                0 TO        0

L4             0 SEA SSS SAM L2 NOT L1

=>
Uploading C:\Program Files\Stnexp\Queries\10823987 (b).str
```



```
chain nodes :
14  15  17  27  28  39  40  51  52  60  61  62  76  77  91  92  106  107  121  122
128
ring nodes :
1  2  3  4  5  6  7  8  9  18  19  20  21  22  23  24  25  26  30  31  32  33  34
35  36  37  38  42  43  44  45  46  47  48  49  50  67  68  69  70  71  72  73  74
75  82  83  84  85  86  87  88  89  90  97  98  99  100  101  102  103  104  105
112  113  114  115  116  117  118  119  120
```

ring/chain nodes :
16 29 41 53 78 93 108 123
chain bonds :
7-14 14-15 15-16 17-60 17-128 24-27 27-28 28-29 36-39 39-40 40-41 48-51
51-52 52-53 60-61 61-62 74-76 76-77 77-78 89-91 91-92 92-93 104-106
106-107 107-108 119-121 121-122 122-123
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 18-19 18-23 19-20 20-21 21-22
22-23 22-24 23-26 24-25 25-26 30-31 30-35 31-32 32-33 33-34 34-35 34-36
35-38 36-37 37-38 42-43 42-47 43-44 44-45 45-46 46-47 46-48 47-50 48-49
49-50 67-68 67-72 68-69 69-70 70-71 71-72 71-73 72-75 73-74 74-75 82-83
82-87 83-84 84-85 85-86 86-87 86-88 87-90 88-89 89-90 97-98 97-102
98-99 99-100 100-101 101-102 101-103 102-105 103-104 104-105 112-113
112-117 113-114 114-115 115-116 116-117 116-118 117-120 118-119 119-120
exact/norm bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 7-14 8-9 14-15 15-16 17-60 17-128
18-19 18-23 19-20 20-21 21-22 22-23 22-24 23-26 24-25 24-27 25-26 27-28
28-29 30-31 30-35 31-32 32-33 33-34 34-35 34-36 35-38 36-37 36-39 37-38
39-40 40-41 42-43 42-47 43-44 44-45 45-46 46-47 46-48 47-50 48-49 48-51
49-50 51-52 52-53 60-61 61-62 67-68 67-72 68-69 69-70 70-71 71-72 71-73
72-75 73-74 74-75 74-76 76-77 77-78 82-83 82-87 83-84 84-85 85-86 86-87
86-88 87-90 88-89 89-90 89-91 91-92 92-93 97-98 97-102 98-99 99-100
100-101 101-102 101-103 102-105 103-104 104-105 104-106 106-107 107-108
112-113 112-117 113-114 114-115 115-116 116-117 116-118 117-120 118-119
119-120 119-121 121-122 122-123

G1:C,N

G2:[*1],[*2],[*3],[*4],[*5],[*6],[*7],[*8]

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 14:CLASS
15:CLASS 16:CLASS 17:CLASS 18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom
24:Atom 25:Atom 26:Atom 27:CLASS 28:CLASS 29:CLASS 30:Atom 31:Atom 32:Atom
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69:Atom 70:Atom 71:Atom 72:Atom 73:Atom 74:Atom 75:Atom 76:CLASS 77:CLASS
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101:Atom 102:Atom 103:Atom 104:Atom 105:Atom 106:CLASS 107:CLASS 108:CLASS
112:Atom 113:Atom 114:Atom 115:Atom 116:Atom 117:Atom 118:Atom 119:Atom
120:Atom 121:CLASS 122:CLASS 123:CLASS 128:CLASS

L5 STRUCTURE UPLOADED

=> d l5
L5 HAS NO ANSWERS
L5 STR
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

```
=> s l5 sss sam
SAMPLE SEARCH INITIATED 17:51:59 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -     12 TO ITERATE

100.0% PROCESSED      12 ITERATIONS                                0 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:           33 TO      447
PROJECTED ANSWERS:               0 TO        0

L6            0 SEA SSS SAM L5

=> s l5 sss ful
FULL SEARCH INITIATED 17:52:09 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -    311 TO ITERATE

100.0% PROCESSED     311 ITERATIONS                                7 ANSWERS
SEARCH TIME: 00.00.01

L7            7 SEA SSS FUL L5

=> => s l7
L8            1 L7

=> d l8 bib,ab,hitstr
```

L8 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN
AN 2000:161121 CAPLUS
DN 132:207763
TI Preparation of benzopyran, tetrahydroquinoline, pyrano[2,3-b]pyridine, and indan derivatives as potassium channel inhibitors
IN Lloyd, John; Finlay, Heather J.; Vaccaro, Wayne; Atwal, Karnail; Gross, Michael F.; Spear, Kerry L.
PA Bristol-Myers Squibb Company, USA
SO PCT Int. Appl., 210 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2000012077	A1	20000309	WO 1999-US18599	19990816
	W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, GM, KE, LS, MW, SD, SL, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
	CA 2341678	AA	20000309	CA 1999-2341678	19990816
	AU 9956753	A1	20000321	AU 1999-56753	19990816
	AU 754204	B2	20021107		
	EP 1109544	A1	20010627	EP 1999-943714	19990816
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
	JP 2002523451	T2	20020730	JP 2000-567195	19990816
	US 6150356	A	20001121	US 1999-375955	19990817
	US 6511977	B1	20030128	US 2000-670285	20000925
	US 2004058931	A1	20040325	US 2002-295574	20021115
	US 2004067944	A1	20040408	US 2002-295404	20021115
	US 6784189	B2	20040831		
PRAI	US 1998-98709P	P	19980901		
	WO 1999-US18599	W	19990816		
	US 1999-375955	A3	19990817		
	US 2000-670285	A3	20000925		

OS MARPAT 132:207763
AB The title compds. (I) [wherein A, B, and D = independently CH or N; R = H, (aryl)alkyl, alkenyl, aryl, (hetero)cycloalkyl, or cycloalkylalkyl; R1 = (aryl)alkyl, aryl, alkenyl, heterocyclo, NR5-heterocyclo, (hetero)cycloalkyl, cycloalkylalkyl, or (un)substituted amino; or R and R1 taken together with the N-S atoms = a 5- to 8-membered ring; R2 = H, (aryl)alkyl, acyl, carboxymethyl, carbamoylmethyl, etc.; R3 and R4 = independently = H, (aryl)alkyl, cycloalkyl, or R3 and R4 taken together with the C to which they are attached form a 5- to 8-membered ring; R5 = H, (aryl)alkyl, alkenyl, aryl, or cycloalkyl(alkyl); X1 = (CR3R4)n, O, NR5, S, S(O), SO2, -OCR3R4-, -NR5CR3R4-, -SCR3R4-, -S(O)CR3R4-, or -SO2CR3R4-; n = 1-3; X2 = single bond, NR5, or O; Q = substituted NHCH:NCN, acyl, (un)substituted sulfamoyl, or substituted heterocyclo] were prepd by solution phase or solid phase synthesis as antiarrhythmics. For example, II was formed in a 3-step sequence involving: (1) sulfonylation of (trans)-4-amino-3,4-dihydro-2,2-dimethyl-6-cyano-2H-benzopyran with 4-ethylbenzenesulfonyl chloride (85%), (2) hydrolysis of

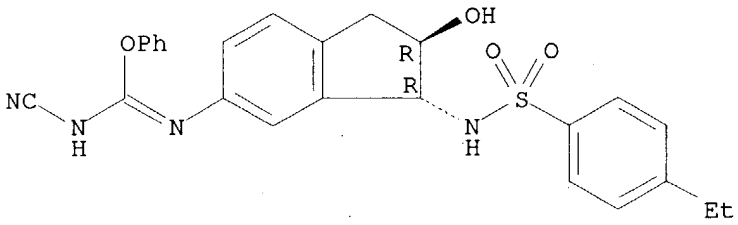
the nitrile to the carboxylic acid using aqueous Na2O2 (33%), and (3) amidation with 1,2,3,4-tetrahydro-1-naphthylamine (51%). I block the delayed rectifier voltage-gated K+ channel (IKur) and are therefore useful in the prevention and treatment of cardiac arrhythmia (no data).

IT **260402-16-2P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(intermediate; preparation of arylsulfamido benzopyran, tetrahydroquinoline, pyrano[2,3-b]pyridine, and indan derivs. by solution phase or solid phase synthesis as potassium channel inhibitors for the treatment of arrhythmia)

RN 260402-16-2 CAPLUS

CN Carbamimidic acid, N-cyano-N'-[(2R,3R)-3-[[(4-ethylphenyl)sulfonyl]amino]-2,3-dihydro-2-hydroxy-1H-inden-5-yl]-, phenyl ester, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

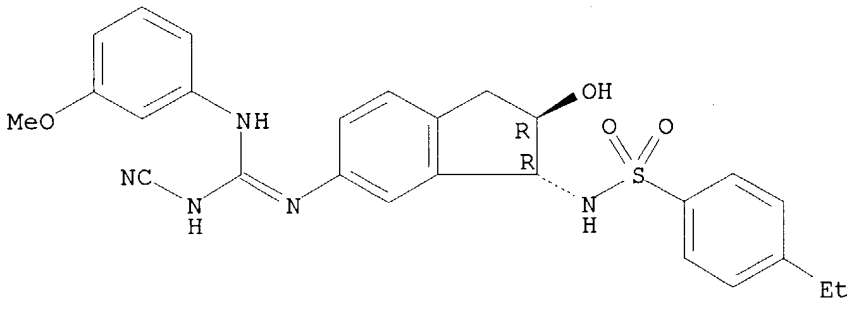


IT **260399-04-0P 260399-05-1P 260399-06-2P**
260399-07-3P 260399-08-4P 260399-09-5P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(target compound; preparation of arylsulfamido benzopyran, tetrahydroquinoline, pyrano[2,3-b]pyridine, and indan derivs. by solution phase or solid phase synthesis as potassium channel inhibitors for the treatment of arrhythmia)

RN 260399-04-0 CAPLUS

CN Benzenesulfonamide, N-[(1R,2R)-6-[[(cyanoamino)[(3-methoxyphenyl)amino]methylene]amino]-2,3-dihydro-2-hydroxy-1H-inden-1-yl]-4-ethyl-, rel- (9CI) (CA INDEX NAME)

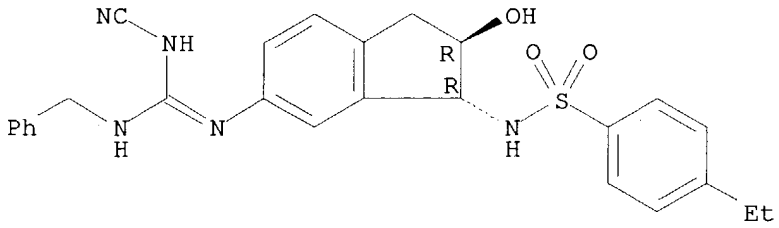
Relative stereochemistry.



RN 260399-05-1 CAPLUS

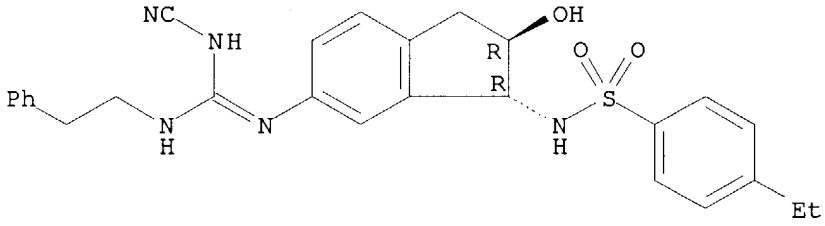
CN Benzenesulfonamide, N-[(1R,2R)-6-[[(cyanoamino)[(phenylmethyl)amino]methylene]amino]-2,3-dihydro-2-hydroxy-1H-inden-1-yl]-4-ethyl-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



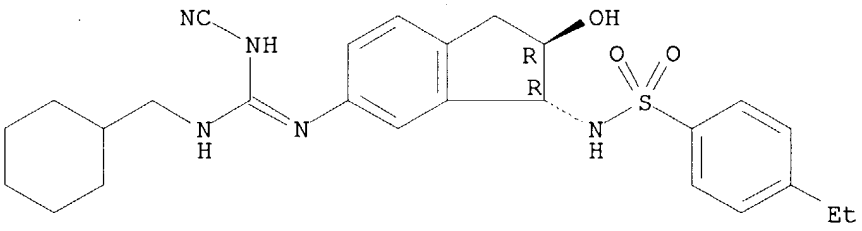
RN 260399-06-2 CAPLUS
CN Benzenesulfonamide, N-[(1R,2R)-6-[[(cyanoamino)[(2-phenylethyl)amino]methylene]amino]-2,3-dihydro-2-hydroxy-1H-inden-1-yl]-4-ethyl-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



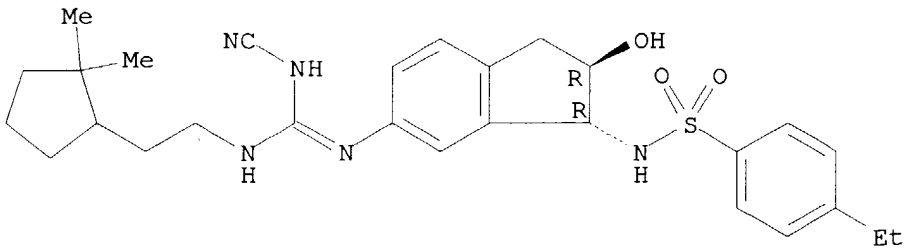
RN 260399-07-3 CAPLUS
CN Benzenesulfonamide, N-[(1R,2R)-6-[[(cyanoamino)[(cyclohexylmethyl)amino]methylene]amino]-2,3-dihydro-2-hydroxy-1H-inden-1-yl]-4-ethyl-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



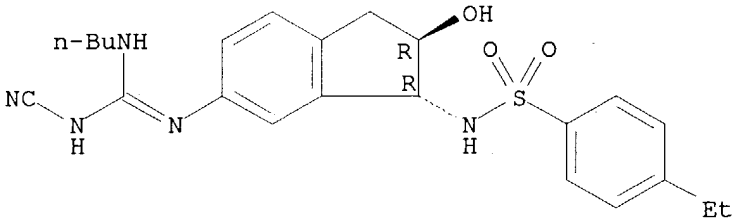
RN 260399-08-4 CAPLUS
CN Benzenesulfonamide, N-[(1R,2R)-6-[[(cyanoamino)[[2-(2,2-dimethylcyclopentyl)ethyl]amino]methylene]amino]-2,3-dihydro-2-hydroxy-1H-inden-1-yl]-4-ethyl-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RN 260399-09-5 CAPLUS
CN Benzenesulfonamide, N-[(1R,2R)-6-[[(butylamino)(cyanoamino)methylene]amino]-2,3-dihydro-2-hydroxy-1H-inden-1-yl]-4-ethyl-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.



RE.CNT 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

```
=> => d his

     (FILE 'HOME' ENTERED AT 17:49:05 ON 23 SEP 2004)

     FILE 'REGISTRY' ENTERED AT 17:49:10 ON 23 SEP 2004
L1             SCREEN 2016 OR  2026 OR  2039 OR  2040 OR  2045 OR  2047
L2             STRUCTURE UPLOADED
L3             QUE L2 NOT L1
L4           0 S L3 SSS SAM
L5             STRUCTURE UPLOADED
L6           0 S L5 SSS SAM
L7           7 S L5 SSS FUL

     FILE 'CAPLUS' ENTERED AT 17:52:16 ON 23 SEP 2004
L8           1 S L7

     FILE 'CAOLD' ENTERED AT 17:52:44 ON 23 SEP 2004

=> s l7
L9           0 L7

=> log y
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                      ENTRY    SESSION
FULL ESTIMATED COST                                    0.42     162.93

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)          SINCE FILE      TOTAL
                                                      ENTRY    SESSION
CA SUBSCRIBER PRICE                                    0.00      -0.70

STN INTERNATIONAL LOGOFF AT 17:52:55 ON 23 SEP 2004
```